## AMENDMENTS TO THE CLAIMS

Please amend the claims so that they read as follows:

 (Previously Presented) A tetrahydro-naphthalene derivative of the formula (I), its tautomeric or stereoisomeric form, or a salt thereof:

$$\begin{array}{c|c} & & & & \\ & & & & \\ & & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ &$$

wherein

n represents an integer of 0 to 6;

R<sub>1</sub> represents hydrogen or C<sub>1-6</sub> alkyl;

 $R_2$  and  $R_3$  together with the nitrogen atom to which they are attached, form a 5-7 membered saturated heterocyclic ring optionally interrupted by one or two atoms selected from the group consisting of oxygen and nitrogen,

wherein said saturated heterocyclic ring has one or more substituents selected from the group consisting of halogen, benzyl, hydroxy, carboxy, amino, oxo, aminocarbonyl, C<sub>1-6</sub>

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alkoxycarbonyl, and  $C_{1-6}$  alkyl optionally substituted by hydroxy, carboxy,  $C_{1-6}$  alkoxy, or  $C_{1-6}$  alkoxycarbonyl,

or

 $R_2$  represents  $C_{2.6}$  alkenyl,  $C_{2.6}$  alkynyl, or  $C_{1.6}$  alkyl substituted by amino, hydroxy,  $C_{1.6}$  alkylamino, or  $di(C_{1.6}$  alkyl)amino;

 $R_3$  represents hydrogen,  $C_{2.6}$  alkenyl,  $C_{2.6}$  alkynyl, or  $C_{1.6}$  alkyl optionally substituted by amino, hydroxy,  $C_{1.6}$  alkylamino, or di( $C_{1.6}$  alkyl)amino; and

 $R_4$  represents hydrogen, halogen,  $C_{1.6}$  alkylthio,  $C_{1.6}$  alkyl optionally substituted by mono-, di-, or tri-halogen, or  $C_{1.6}$  alkoxy optionally substituted by mono-, di-, or tri-halogen.

2. (Previously Presented) The tetrahydro-naphthalene derivative of the formula (I), its tautomeric or stereoisomeric form, or a salt thereof as claimed in claim 1,

wherein

n represents an integer of 0 or 1;

R<sub>1</sub> represents hydrogen;

 $R_2$  and  $R_3$  together with the nitrogen atom to which they are attached, form a 5-7 membered saturated heterocyclic ring optionally interrupted by one or two atoms selected from the group consisting of oxygen and nitrogen,

wherein said saturated heterocyclic ring has one or more substituents selected from the group consisting of benzyl, hydroxy, carboxy, oxo, aminocarbonyl,  $C_{1-6}$  alkoxycarbonyl, and  $C_{1-6}$  alkyl optionally substituted by hydroxy,  $C_{1-6}$  alkoxy, or  $C_{1-6}$  alkoxycarbonyl,

or

R<sub>2</sub> represents C<sub>1-6</sub> alkyl substituted by hydroxy, amino, C<sub>1-6</sub> alkylamino, or di(C<sub>1-6</sub> alkyl)amino;

 $R_3$  represents hydrogen,  $C_{1.6}$  alkyl optionally substituted by hydroxy, amino,  $C_{1.6}$  alkylamino, or  $di(C_{1.6}$  alkyl)amino; and

 $R_4$  represents hydrogen, halogen,  $C_{1-6}$  alkyl optionally substituted by mono-, di-, or tri-halogen, or  $C_{1-6}$  alkoxy optionally substituted by mono-, di-, or tri-halogen.

3. (Previously Presented) The tetrahydro-naphthalene derivative of the formula (I), its tautomeric or stereoisomeric form, or a salt thereof as claimed in claim 1, wherein

n represents an integer of 0 or 1;

R<sub>1</sub> represents hydrogen;

 $R_2$  and  $R_3$  together with the nitrogen atom to which they are attached, form a pyrrolidinyl optionally substituted by oxo, piperidinyl optionally substituted by hydroxy, carboxy, aminocarbonyl,  $C_{1-6}$  alkoxycarbonyl, or  $C_{1-6}$  alkyl optionally substituted by hydroxy, piperazinyl optionally substituted by benzyl, homopiperidinyl, or morpholinyl,

or

 $R_2$  represents  $C_{1-6}$  alkyl substituted by hydroxy, or  $di(C_{1-6}$  alkyl)amino;  $R_3$  represents hydrogen, or  $C_{1-6}$  alkyl; and  $R_4$  represents hydrogen, fluoro, chloro, bromo,  $C_{1-6}$  alkyl optionally substituted by mono-, di-, or tri-halogen, or  $C_{1-6}$  alkoxy.

4. (Previously Presented) A tetrahydro-naphthalene derivative, its tautomeric or stereoisomeric form, or a salt thereof selected from the group consisting of:

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N-(7-hydroxy-5,6,7,8-tetrahydronaphthalen-1-yl)-N'-[3-piperidin-1-yl-4-(trifluoromethyl)benzyl]urea;

N-(7-hydroxy-5,6,7,8-tetrahydronaphthalen-1-yl)-N'-[4-pyrrolidin-1-yl-3-(trifluoromethyl)benzyl]urea;

N-(7-hydroxy-5,6,7,8-tetrahydronaphthalen-1-yl)-N'-[3-pyrrolidin-1-yl-4-(trifluoromethyl)benzyl]urea;

 $\label{eq:n-l-yl-3-(trifluoromethyl)benzyl]-N'-(7-hydroxy-5,6,7,8-tetrahydronaphthalen-1-yl)urea;$ 

N-[3-azepan-1-yl-4-(trifluoromethyl)benzyl]-N'-(7-hydroxy-5,6,7,8-tetrahydronaphthalen-1-yl)urea;

N-(3-bromo-4-piperidin-1-ylbenzyl)-N'-(7-hydroxy-5,6,7,8-tetrahydronaphthalen-1-yl)urea;

 $\label{eq:n-index} N-[(7R)-7-hydroxy-5,6,7,8-tetrahydronaphthalen-1-yl]-N'-[3-pyrrolidin-1-yl-4-(trifluoromethyl)benzyl]urea;$ 

N-[(7S)-7-hydroxy-5,6,7,8-tetrahydronaphthalen-1-yl]-N'-[3-pyrrolidin-1-yl-4-(trifluoromethyl)benzyl]urea;

 $\label{eq:continuous} N-(7-hydroxy-5,6,7,8-tetrahydronaphthalen-1-yl)-N'-[4-piperidin-1-yl-3-(trifluoromethyl)benzyl]urea;$ 

ethyl 1-[5-[({[(7-hydroxy-5,6,7,8-tetrahydronaphthalen-1-yl)amino]carbonyl}amino)-methyl]-2-(trifluoromethyl)phenyl]piperidine-4-carboxylate; and

N-[(7R)-7-hydroxy-5,6,7,8-tetra hydron aphthalen-1-yl]-N'-[3-morpholin-4-yl-4-(trifluor omethyl) benzyl]urea.

5. (Previously Presented) A pharmaceutical composition comprising a tetrahydro-naphthalene derivative of the formula (I), its tautomeric or stereoisomeric form, or a physiologically acceptable salt thereof as claimed in claim 1 as an active ingredient, and a pharmaceutically acceptable excipient.

Claims 6 - 26, (canceled)

- 27. (Previously Presented) The tetrahydro-napthalene derivative of claim 4, its tautomeric or stereoisomeric form, or a salt thereof wherein the compound is N-[(7R)-7-hydroxy-5,6,7,8-tetrahydronaphthalen-1-yll-N-[3-pyrrolidin-1-yl-4-(trifluoromethyl)benzyl]urea.
- 28. (Previously Presented) The tetrahydro-napthalene derivative of claim 4, its tautomeric or stereoisomeric form, or a salt thereof wherein the compound is N-[(7R)-7-hydroxy-5,6,7,8-tetrahydronaphthalen-1-yl]-N-[3-morpholin-4-yl-4-(trifluoromethyl)benzyl]urea.

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